



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:38 PM GMT

PDB ID : 1GGE  
Title : CRYSTAL STRUCTURE OF CATALASE HP11 FROM ESCHERICHIA COLI, NATIVE STRUCTURE AT 1.9 Å RESOLUTION.  
Authors : Melik-Adamyan, W.R.; Bravo, J.; Carpena, X.; Switala, J.; Mate, M.J.; Fita, I.; Loewen, P.C.  
Deposited on : 2000-08-16  
Resolution : 1.89 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

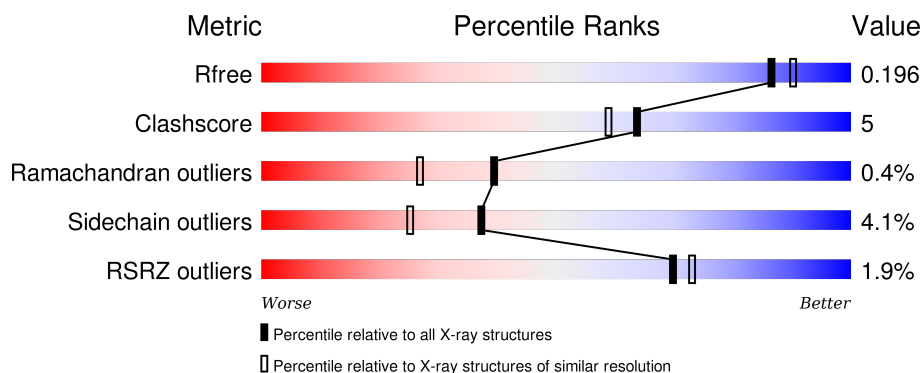
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.89 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	753	<div> <div>2%</div> <div>84%</div> <div>11%</div> <div>• •</div> </div>
1	B	753	<div> <div>2%</div> <div>82%</div> <div>13%</div> <div>• •</div> </div>
1	C	753	<div> <div>2%</div> <div>82%</div> <div>13%</div> <div>• •</div> </div>
1	D	753	<div> <div>%</div> <div>81%</div> <div>14%</div> <div>• •</div> </div>

## 2 Entry composition i

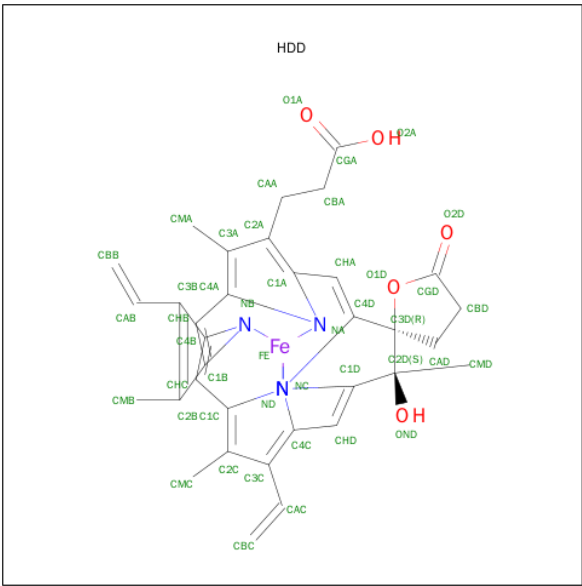
There are 3 unique types of molecules in this entry. The entry contains 26368 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called PROTEIN (CATALASE HP1I).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	727	Total	C	N	O	S	0	0	0
			5746	3647	1005	1082	12			
1	B	727	Total	C	N	O	S	0	0	0
			5746	3647	1005	1082	12			
1	C	727	Total	C	N	O	S	0	0	0
			5746	3647	1005	1082	12			
1	D	727	Total	C	N	O	S	0	0	0
			5746	3647	1005	1082	12			

- Molecule 2 is CIS-HEME D HYDROXYCHLORIN GAMMA-SPIROLACTONE (three-letter code: HDD) (formula: C<sub>34</sub>H<sub>32</sub>FeN<sub>4</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		
2	B	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		
2	D	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		

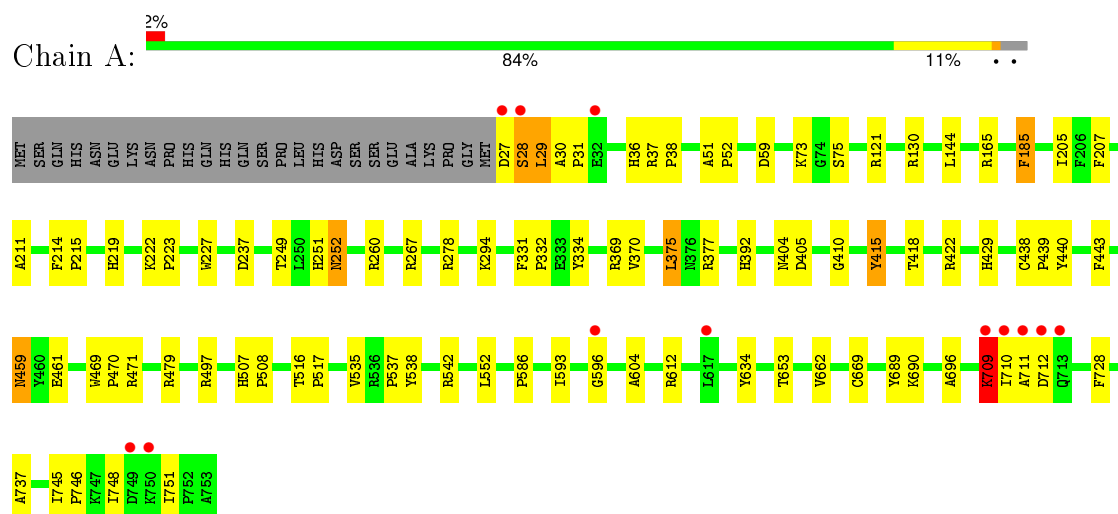
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	843	Total	O	0	0
			843	843		
3	B	735	Total	O	0	0
			735	735		
3	C	791	Total	O	0	0
			791	791		
3	D	839	Total	O	0	0
			839	839		

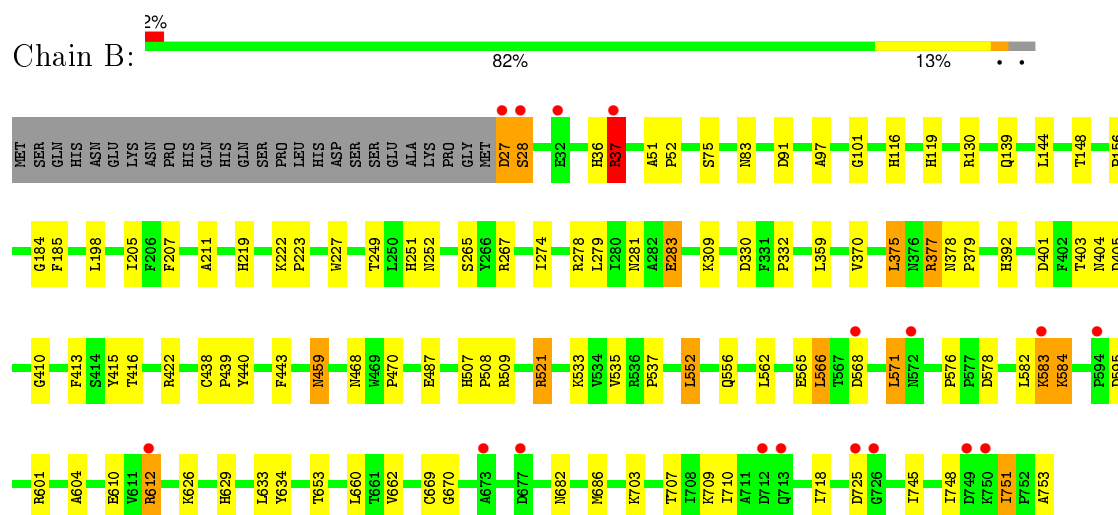
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

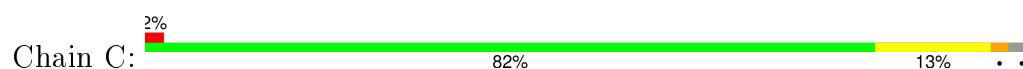
#### • Molecule 1: PROTEIN (CATALASE HP1I)

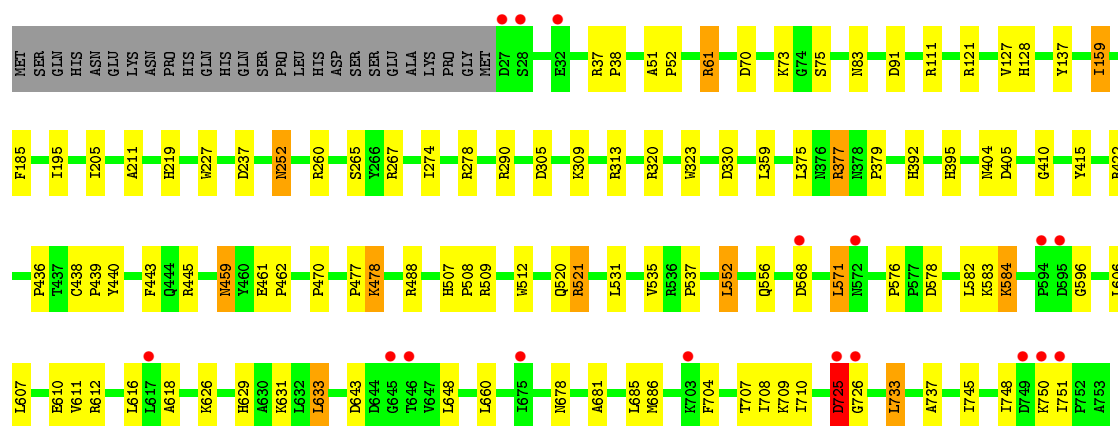


#### • Molecule 1: PROTEIN (CATALASE HP1I)

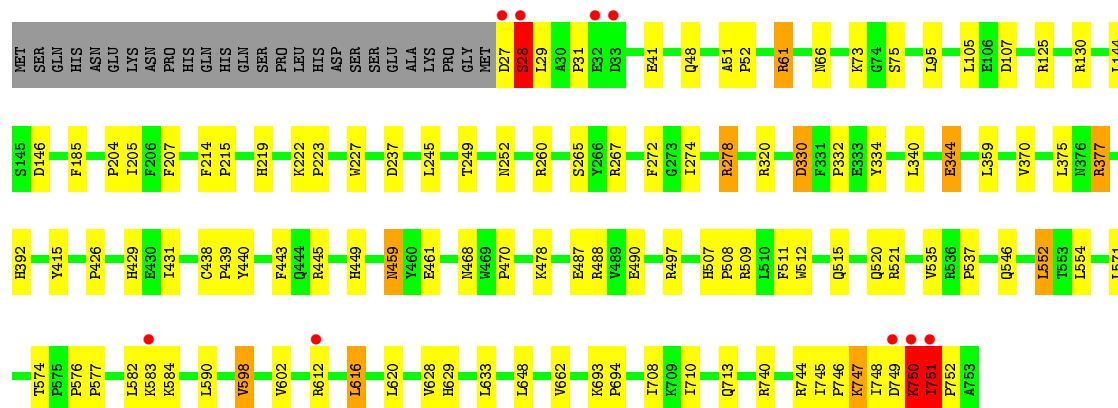
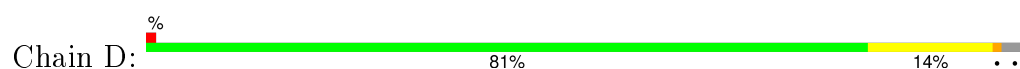


#### • Molecule 1: PROTEIN (CATALASE HP1I)





• Molecule 1: PROTEIN (CATALASE HP11)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.04Å 132.34Å 121.20Å 90.00° 109.63° 90.00°	Depositor
Resolution (Å)	87.60 – 1.89 17.96 – 1.89	Depositor EDS
% Data completeness (in resolution range)	(Not available) (87.60-1.89) 97.7 (17.96-1.89)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.19 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.163 , 0.202 0.155 , 0.196	Depositor DCC
$R_{free}$ test set	10774 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	9.5	Xtriage
Anisotropy	0.249	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 59.2	EDS
Estimated twinning fraction	0.020 for h,-k,-h-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 214389 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	26368	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HDD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.42	0/5902	1.07	20/8024 (0.2%)
1	B	0.42	0/5902	1.07	9/8024 (0.1%)
1	C	0.43	0/5902	1.07	17/8024 (0.2%)
1	D	0.42	0/5902	1.11	22/8024 (0.3%)
All	All	0.42	0/23608	1.08	68/32096 (0.2%)

There are no bond length outliers.

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	61	ARG	NE-CZ-NH2	12.37	126.48	120.30
1	D	612	ARG	CD-NE-CZ	11.02	139.03	123.60
1	A	121	ARG	NE-CZ-NH2	-8.92	115.84	120.30
1	B	28	SER	N-CA-CB	8.73	123.60	110.50
1	D	61	ARG	NE-CZ-NH1	-8.72	115.94	120.30
1	B	521	ARG	NE-CZ-NH1	8.32	124.46	120.30
1	A	278	ARG	NE-CZ-NH1	-8.20	116.20	120.30
1	C	121	ARG	NE-CZ-NH1	8.17	124.38	120.30
1	D	125	ARG	NE-CZ-NH2	-8.09	116.25	120.30
1	D	377	ARG	NE-CZ-NH2	-8.07	116.26	120.30
1	B	422	ARG	NE-CZ-NH1	7.96	124.28	120.30
1	D	449	HIS	CA-CB-CG	7.84	126.93	113.60
1	D	740	ARG	NE-CZ-NH1	7.77	124.19	120.30
1	A	121	ARG	NE-CZ-NH1	7.66	124.13	120.30
1	C	726	GLY	N-CA-C	-7.34	94.75	113.10
1	A	130	ARG	NE-CZ-NH2	-7.24	116.68	120.30
1	A	377	ARG	NE-CZ-NH2	-7.02	116.79	120.30
1	D	740	ARG	NE-CZ-NH2	-6.96	116.82	120.30
1	D	130	ARG	NE-CZ-NH2	-6.69	116.96	120.30
1	B	601	ARG	NE-CZ-NH2	-6.63	116.98	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	61	ARG	CD-NE-CZ	6.51	132.71	123.60
1	D	445	ARG	NE-CZ-NH2	-6.50	117.05	120.30
1	C	725	ASP	N-CA-CB	6.49	122.28	110.60
1	D	521	ARG	NE-CZ-NH2	6.41	123.50	120.30
1	D	612	ARG	CG-CD-NE	6.39	125.21	111.80
1	D	320	ARG	NE-CZ-NH2	-6.36	117.12	120.30
1	C	61	ARG	NE-CZ-NH1	-6.34	117.13	120.30
1	C	111	ARG	NE-CZ-NH1	-6.27	117.16	120.30
1	D	125	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	C	70	ASP	CB-CG-OD2	6.25	123.92	118.30
1	C	111	ARG	NE-CZ-NH2	6.14	123.37	120.30
1	A	497	ARG	CD-NE-CZ	6.10	132.14	123.60
1	B	377	ARG	NE-CZ-NH1	-6.07	117.27	120.30
1	A	130	ARG	NE-CZ-NH1	6.05	123.32	120.30
1	A	267	ARG	NE-CZ-NH2	-5.97	117.31	120.30
1	B	37	ARG	CD-NE-CZ	5.96	131.95	123.60
1	A	422	ARG	NE-CZ-NH2	5.93	123.27	120.30
1	D	278	ARG	NE-CZ-NH1	-5.93	117.34	120.30
1	B	130	ARG	NE-CZ-NH1	5.89	123.24	120.30
1	A	260	ARG	NE-CZ-NH2	-5.83	117.39	120.30
1	A	369	ARG	NE-CZ-NH2	5.80	123.20	120.30
1	A	542	ARG	NE-CZ-NH2	-5.69	117.45	120.30
1	C	278	ARG	NE-CZ-NH1	-5.67	117.47	120.30
1	D	334	TYR	CB-CG-CD2	-5.65	117.61	121.00
1	C	265	SER	N-CA-CB	-5.65	102.03	110.50
1	C	320	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	C	422	ARG	NE-CZ-NH2	5.53	123.06	120.30
1	A	278	ARG	NE-CZ-NH2	5.50	123.05	120.30
1	B	401	ASP	CB-CG-OD2	5.50	123.25	118.30
1	A	479	ARG	CD-NE-CZ	5.42	131.19	123.60
1	A	334	TYR	CB-CG-CD2	-5.42	117.75	121.00
1	C	290	ARG	NE-CZ-NH1	-5.39	117.61	120.30
1	D	497	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	D	445	ARG	NE-CZ-NH1	5.33	122.96	120.30
1	C	445	ARG	NE-CZ-NH1	5.31	122.95	120.30
1	C	260	ARG	NE-CZ-NH2	-5.25	117.67	120.30
1	B	403	THR	N-CA-CB	5.24	120.26	110.30
1	A	185	PHE	CB-CG-CD2	-5.18	117.17	120.80
1	C	313	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	C	320	ARG	CD-NE-CZ	5.16	130.82	123.60
1	D	344	GLU	CA-CB-CG	5.14	124.71	113.40
1	C	377	ARG	NE-CZ-NH1	-5.10	117.75	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	488	ARG	CD-NE-CZ	5.07	130.70	123.60
1	A	185	PHE	CB-CG-CD1	5.06	124.34	120.80
1	A	59	ASP	CB-CG-OD1	5.05	122.85	118.30
1	A	369	ARG	NE-CZ-NH1	-5.03	117.78	120.30
1	D	330	ASP	CB-CG-OD2	-5.03	113.77	118.30
1	A	165	ARG	NE-CZ-NH2	-5.00	117.80	120.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5746	0	5578	51	0
1	B	5746	0	5578	69	0
1	C	5746	0	5578	53	0
1	D	5746	0	5578	68	0
2	A	44	0	31	6	0
2	B	44	0	31	3	0
2	C	44	0	31	2	0
2	D	44	0	31	1	0
3	A	843	0	0	3	0
3	B	735	0	0	8	0
3	C	791	0	0	8	0
3	D	839	0	0	9	0
All	All	26368	0	22436	231	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (231) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:760:HDD:C3D	2:A:760:HDD:CAD	1.76	1.57
1:A:392:HIS:ND1	1:A:415:TYR:HB2	1.09	1.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:392:HIS:ND1	1:C:415:TYR:HB2	1.08	1.39
1:D:392:HIS:ND1	1:D:415:TYR:HB2	1.10	1.38
1:B:392:HIS:ND1	1:B:415:TYR:HB2	1.08	1.35
1:D:392:HIS:CE1	1:D:415:TYR:HB2	2.11	0.83
2:A:760:HDD:C4D	2:A:760:HDD:CAD	2.57	0.82
1:B:392:HIS:CE1	1:B:415:TYR:HB2	2.15	0.79
1:D:750:LYS:HD3	1:D:751:ILE:H	1.48	0.78
1:B:612:ARG:HD3	1:B:670:GLY:HA2	1.65	0.77
1:A:392:HIS:CE1	1:A:415:TYR:HB2	2.14	0.76
1:B:583:LYS:NZ	1:B:583:LYS:H	1.83	0.74
1:C:552:LEU:HD21	1:C:571:LEU:HD12	1.70	0.74
1:A:392:HIS:CG	1:A:415:TYR:HB2	2.17	0.74
2:C:760:HDD:HBC1	2:C:760:HDD:HMC1	1.72	0.72
1:C:392:HIS:CE1	1:C:415:TYR:HB2	2.14	0.70
1:C:392:HIS:CG	1:C:415:TYR:HB2	2.19	0.69
1:D:267:ARG:HG3	3:D:1253:HOH:O	1.92	0.69
1:C:748:ILE:O	1:C:751:ILE:HG22	1.92	0.69
1:B:552:LEU:HD21	1:B:571:LEU:HD12	1.74	0.69
1:D:392:HIS:CG	1:D:415:TYR:HB2	2.21	0.69
1:B:392:HIS:CG	1:B:415:TYR:HB2	2.20	0.69
1:B:468:ASN:HD22	1:D:27:ASP:N	1.92	0.68
1:C:708:ILE:HG13	1:C:710:ILE:HG12	1.76	0.67
1:D:27:ASP:O	1:D:28:SER:HB2	1.95	0.66
1:B:552:LEU:HD22	1:B:556:GLN:HG3	1.78	0.66
1:B:281:ASN:OD1	1:B:283:GLU:HG3	1.96	0.66
1:B:583:LYS:O	1:B:584:LYS:HB3	1.96	0.66
1:C:267:ARG:HG3	3:C:1464:HOH:O	1.94	0.66
1:B:267:ARG:HG3	3:B:1236:HOH:O	1.95	0.64
1:D:750:LYS:HD3	1:D:751:ILE:N	2.12	0.64
2:A:760:HDD:C2D	2:A:760:HDD:CAD	2.66	0.62
1:B:359:LEU:H	1:B:507:HIS:HD2	1.47	0.61
1:A:690:LYS:HG3	1:A:751:ILE:HD11	1.82	0.61
1:A:748:ILE:O	1:A:751:ILE:HG13	2.00	0.61
1:C:274:ILE:HD12	2:C:760:HDD:HMB1	1.83	0.61
1:D:51:ALA:HB1	1:D:52:PRO:HD2	1.83	0.61
1:D:144:LEU:HD11	1:D:370:VAL:HG13	1.83	0.60
1:D:61:ARG:HH11	1:D:66:ASN:HA	1.66	0.60
1:A:709:LYS:HA	1:A:709:LYS:HE3	1.86	0.58
1:D:27:ASP:HB2	1:D:29:LEU:HD21	1.85	0.57
1:B:710:ILE:HD13	1:B:718:ILE:HG13	1.84	0.57
1:D:583:LYS:O	1:D:584:LYS:HB3	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:578:ASP:HB3	1:B:582:LEU:O	2.05	0.56
1:A:751:ILE:O	1:A:751:ILE:HD12	2.04	0.56
1:A:28:SER:CB	1:D:245:LEU:HD22	2.35	0.56
1:D:708:ILE:HG13	1:D:710:ILE:HG12	1.87	0.56
1:A:28:SER:HB2	1:D:245:LEU:HD13	1.87	0.55
1:D:330:ASP:OD1	1:D:629:HIS:HE1	1.89	0.55
1:C:137:TYR:HB2	1:C:159:ILE:CD1	2.37	0.55
1:C:745:ILE:HD13	3:C:1466:HOH:O	2.07	0.55
1:A:294:LYS:NZ	3:A:1120:HOH:O	2.40	0.54
1:D:260:ARG:HD3	1:D:590:LEU:HD21	1.90	0.54
1:B:51:ALA:HB1	1:B:52:PRO:HD2	1.90	0.54
1:B:583:LYS:HZ3	1:B:583:LYS:H	1.52	0.54
1:D:535:VAL:O	1:D:537:PRO:HD3	2.06	0.54
3:A:912:HOH:O	1:C:52:PRO:HG3	2.08	0.54
1:A:30:ALA:HB1	1:A:31:PRO:HD2	1.90	0.54
1:C:535:VAL:O	1:C:537:PRO:HD3	2.07	0.53
1:A:36:HIS:CD2	1:A:36:HIS:H	2.26	0.53
1:B:330:ASP:OD1	1:B:629:HIS:HE1	1.90	0.53
1:C:610:GLU:OE1	1:C:643:ASP:HA	2.07	0.53
1:A:222:LYS:HB3	1:A:223:PRO:HD2	1.90	0.53
1:D:629:HIS:HD2	3:D:1134:HOH:O	1.91	0.53
1:C:477:PRO:HB2	1:C:478:LYS:HD3	1.91	0.53
1:A:459:ASN:ND2	1:B:219:HIS:HB3	2.25	0.52
1:B:37:ARG:HD2	3:B:1429:HOH:O	2.09	0.52
2:B:760:HDD:HBB1	2:B:760:HDD:HMB1	1.91	0.52
1:C:443:PHE:CZ	1:C:470:PRO:HD2	2.45	0.52
1:D:359:LEU:H	1:D:507:HIS:HD2	1.56	0.52
1:A:27:ASP:O	1:A:28:SER:C	2.48	0.52
1:B:27:ASP:HB3	3:B:1359:HOH:O	2.10	0.52
1:C:359:LEU:H	1:C:507:HIS:HD2	1.58	0.52
1:C:583:LYS:O	1:C:584:LYS:HB3	2.10	0.51
1:D:509:ARG:HD2	1:D:576:PRO:HD2	1.93	0.51
1:C:578:ASP:HB2	1:C:582:LEU:O	2.11	0.51
1:D:744:ARG:HA	1:D:747:LYS:HD3	1.91	0.51
1:B:521:ARG:HH21	1:B:745:ILE:HD13	1.74	0.51
3:B:961:HOH:O	1:D:52:PRO:HG3	2.09	0.51
1:B:521:ARG:HD3	3:B:1462:HOH:O	2.11	0.50
1:B:535:VAL:O	1:B:537:PRO:HD3	2.11	0.50
1:D:552:LEU:HD11	1:D:571:LEU:HD23	1.92	0.50
1:A:29:LEU:HB2	3:C:1352:HOH:O	2.10	0.50
1:B:416:THR:HG21	3:D:1400:HOH:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:HIS:HD1	1:B:36:HIS:H	1.60	0.50
1:C:211:ALA:CB	1:C:410:GLY:HA3	2.41	0.50
1:A:418:THR:HG21	2:A:760:HDD:HBD2	1.93	0.50
1:B:443:PHE:CZ	1:B:470:PRO:HD2	2.46	0.50
1:D:602:VAL:HG13	1:D:662:VAL:HA	1.92	0.50
1:B:556:GLN:HG2	1:B:566:LEU:HD23	1.93	0.49
1:D:598:VAL:HG13	1:D:628:VAL:CG2	2.42	0.49
1:A:586:PRO:HB2	1:A:593:ILE:HD11	1.93	0.49
1:B:359:LEU:H	1:B:507:HIS:CD2	2.29	0.49
1:D:616:LEU:HD22	1:D:620:LEU:HG	1.95	0.49
1:D:478:LYS:HD2	3:D:1546:HOH:O	2.12	0.49
1:A:535:VAL:O	1:A:537:PRO:HD3	2.13	0.49
1:D:214:PHE:HB3	1:D:215:PRO:HD3	1.94	0.49
1:D:146:ASP:HB2	3:D:1383:HOH:O	2.13	0.48
1:C:330:ASP:OD1	1:C:629:HIS:HE1	1.95	0.48
1:A:219:HIS:HB3	1:B:459:ASN:ND2	2.28	0.48
1:D:546:GLN:HG3	3:D:1486:HOH:O	2.13	0.48
1:C:626:LYS:HG3	1:C:733:LEU:HG	1.95	0.48
1:D:61:ARG:NH1	1:D:66:ASN:HA	2.29	0.48
1:D:751:ILE:HD13	1:D:752:PRO:HD2	1.96	0.48
1:C:459:ASN:ND2	1:D:219:HIS:HB3	2.29	0.48
1:D:274:ILE:HD12	2:D:760:HDD:HMB1	1.96	0.48
1:B:710:ILE:CD1	1:B:718:ILE:HG13	2.44	0.48
1:B:27:ASP:OD2	1:D:468:ASN:ND2	2.37	0.48
1:B:251:HIS:CE1	1:B:507:HIS:HB3	2.49	0.47
1:B:207:PHE:O	1:B:249:THR:HA	2.14	0.47
1:C:704:PHE:O	1:C:707:THR:HG22	2.14	0.47
1:A:214:PHE:HB3	1:A:215:PRO:HD3	1.97	0.47
1:A:443:PHE:CZ	1:A:470:PRO:HD2	2.49	0.47
1:A:211:ALA:CB	1:A:410:GLY:HA3	2.45	0.47
1:C:219:HIS:HB3	1:D:459:ASN:ND2	2.30	0.47
1:B:97:ALA:O	1:B:101:GLY:HA3	2.14	0.47
1:B:278:ARG:HH22	1:B:487:GLU:CD	2.18	0.46
1:A:404:ASN:O	1:A:405:ASP:C	2.54	0.46
1:C:629:HIS:HD2	3:C:1067:HOH:O	1.99	0.46
1:B:144:LEU:HD11	1:B:370:VAL:HG13	1.97	0.46
1:D:431:ILE:HD11	3:D:1575:HOH:O	2.16	0.46
1:B:116:HIS:CD2	1:D:426:PRO:HB2	2.51	0.46
1:D:207:PHE:O	1:D:249:THR:HA	2.15	0.46
1:C:751:ILE:HB	3:C:1343:HOH:O	2.15	0.46
1:B:533:LYS:HE2	3:B:1471:HOH:O	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:HIS:ND1	1:A:415:TYR:CG	2.80	0.46
1:B:604:ALA:HB2	1:B:662:VAL:HG11	1.97	0.46
1:A:612:ARG:HG3	3:A:1460:HOH:O	2.15	0.46
1:C:610:GLU:O	1:C:610:GLU:HG3	2.15	0.46
1:B:309:LYS:HD2	1:B:660:LEU:HD11	1.98	0.46
1:B:459:ASN:H	1:B:459:ASN:HD22	1.64	0.45
1:C:38:PRO:HG2	1:C:51:ALA:HB2	1.98	0.45
1:B:413:PHE:HB2	1:D:105:LEU:HD11	1.98	0.45
1:C:359:LEU:H	1:C:507:HIS:CD2	2.35	0.45
1:A:332:PRO:HD2	1:A:375:LEU:O	2.17	0.45
1:D:267:ARG:HG2	1:D:332:PRO:HB3	1.98	0.45
1:C:323:TRP:CZ3	1:C:379:PRO:HD2	2.51	0.45
1:D:438:CYS:HB2	1:D:439:PRO:CD	2.47	0.45
1:C:686:MET:HB3	1:C:751:ILE:HD11	1.99	0.45
1:A:28:SER:HB2	1:D:245:LEU:HD22	1.99	0.45
1:C:521:ARG:HG2	3:C:1466:HOH:O	2.17	0.45
1:C:618:ALA:HB1	1:C:725:ASP:HB2	1.99	0.45
1:C:73:LYS:HD2	3:C:1424:HOH:O	2.17	0.44
1:B:629:HIS:HD2	3:B:1029:HOH:O	1.99	0.44
1:D:574:THR:HG22	3:D:1182:HOH:O	2.16	0.44
1:C:392:HIS:HB3	1:C:395:HIS:CE1	2.53	0.44
1:A:51:ALA:HB1	1:A:52:PRO:HD2	1.99	0.44
1:D:443:PHE:CZ	1:D:470:PRO:HD2	2.53	0.44
1:B:751:ILE:HD12	1:B:753:ALA:HB2	1.99	0.44
1:D:222:LYS:HB3	1:D:223:PRO:HD2	1.99	0.44
1:A:52:PRO:HG3	3:C:1000:HOH:O	2.16	0.44
1:B:438:CYS:HB2	1:B:439:PRO:HD2	1.98	0.44
1:C:686:MET:CB	1:C:751:ILE:HD11	2.47	0.44
1:B:521:ARG:NH2	1:B:745:ILE:HD13	2.32	0.44
1:C:305:ASP:O	1:C:309:LYS:HG3	2.17	0.44
1:A:696:ALA:HB1	1:A:728:PHE:CZ	2.53	0.44
1:A:37:ARG:HA	1:A:38:PRO:HD3	1.83	0.44
1:B:211:ALA:CB	1:B:410:GLY:HA3	2.48	0.44
1:A:612:ARG:HE	1:A:669:CYS:HB3	1.83	0.44
1:B:748:ILE:O	1:B:751:ILE:HG22	2.18	0.43
1:A:745:ILE:N	1:A:746:PRO:HD2	2.33	0.43
1:A:222:LYS:HB3	1:A:223:PRO:CD	2.48	0.43
1:C:438:CYS:HB2	1:C:439:PRO:HD2	2.00	0.43
1:A:207:PHE:O	1:A:249:THR:HA	2.18	0.43
1:C:678:ASN:HB3	1:C:681:ALA:HB3	1.99	0.43
1:C:596:GLY:HA3	1:C:737:ALA:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:689:TYR:OH	1:A:710:ILE:HG21	2.19	0.43
1:C:552:LEU:HD13	1:C:556:GLN:NE2	2.34	0.43
1:C:607:LEU:HD22	1:C:611:VAL:HG21	2.01	0.43
1:D:745:ILE:HB	1:D:746:PRO:HD3	1.99	0.43
1:B:686:MET:HG2	1:B:751:ILE:HD11	2.00	0.43
1:C:127:VAL:O	1:C:128:HIS:HB2	2.19	0.43
1:A:538:TYR:CE2	1:D:31:PRO:HG3	2.54	0.43
1:D:693:LYS:HA	1:D:694:PRO:HD3	1.85	0.43
1:C:252:ASN:HD22	1:C:252:ASN:HA	1.67	0.43
1:D:748:ILE:O	1:D:751:ILE:HG22	2.18	0.43
1:B:274:ILE:HD12	2:B:760:HDD:HMB1	2.00	0.43
1:B:634:TYR:O	1:B:653:THR:HA	2.18	0.43
1:D:95:LEU:HB3	1:D:107:ASP:HB2	1.99	0.43
1:D:577:PRO:HG2	3:D:1579:HOH:O	2.19	0.43
1:C:404:ASN:O	1:C:405:ASP:C	2.55	0.43
1:D:507:HIS:N	1:D:508:PRO:CD	2.82	0.42
1:C:195:ILE:HD11	1:C:436:PRO:HA	2.01	0.42
1:B:83:ASN:HB3	1:D:429:HIS:CG	2.53	0.42
1:C:509:ARG:HD2	1:C:576:PRO:HD2	2.00	0.42
1:B:139:GLN:HG2	1:B:156:PRO:O	2.19	0.42
1:D:507:HIS:N	1:D:508:PRO:HD2	2.34	0.42
1:B:686:MET:SD	1:B:707:THR:HG22	2.59	0.42
1:A:469:TRP:CE3	1:A:471:ARG:HG3	2.54	0.42
1:B:332:PRO:HD2	1:B:375:LEU:O	2.19	0.42
1:D:616:LEU:HD23	1:D:616:LEU:HA	1.93	0.42
1:B:119:HIS:HB2	1:D:426:PRO:HG3	2.02	0.42
1:A:252:ASN:HA	1:A:252:ASN:HD22	1.71	0.42
1:A:596:GLY:HA3	1:A:737:ALA:O	2.19	0.42
1:C:631:LYS:HG3	1:C:633:LEU:HD13	2.02	0.42
1:A:604:ALA:HB2	1:A:662:VAL:HG11	2.01	0.42
1:A:438:CYS:HB2	1:A:439:PRO:CD	2.50	0.42
1:A:634:TYR:O	1:A:653:THR:HA	2.19	0.42
1:D:392:HIS:ND1	1:D:415:TYR:CG	2.83	0.42
1:B:507:HIS:N	1:B:508:PRO:CD	2.82	0.42
1:D:204:PRO:HB3	1:D:272:PHE:CD1	2.55	0.42
1:B:222:LYS:HB3	1:B:223:PRO:HD2	2.02	0.42
1:D:512:TRP:CH2	1:D:520:GLN:HB3	2.55	0.41
1:B:404:ASN:O	1:B:405:ASP:C	2.59	0.41
1:B:509:ARG:HD2	1:B:576:PRO:HD2	2.03	0.41
2:A:760:HDD:HBB1	2:A:760:HDD:HMB1	2.00	0.41
1:D:359:LEU:H	1:D:507:HIS:CD2	2.37	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:709:LYS:HD2	1:B:709:LYS:HA	1.74	0.41
1:B:521:ARG:HH21	1:B:521:ARG:HD2	1.72	0.41
1:B:378:ASN:HB3	1:B:379:PRO:HD2	2.03	0.41
1:A:507:HIS:N	1:A:508:PRO:CD	2.84	0.41
1:A:516:THR:HB	1:A:517:PRO:HD2	2.02	0.41
1:D:27:ASP:HB2	1:D:29:LEU:CD2	2.50	0.41
1:C:507:HIS:N	1:C:508:PRO:CD	2.83	0.41
1:A:331:PHE:HA	1:A:332:PRO:HD3	1.95	0.41
1:D:222:LYS:HB3	1:D:223:PRO:CD	2.51	0.41
1:B:626:LYS:NZ	3:B:1341:HOH:O	2.54	0.41
1:A:429:HIS:CG	1:C:83:ASN:HB3	2.55	0.41
1:C:461:GLU:HA	1:C:462:PRO:C	2.41	0.41
1:A:251:HIS:CE1	1:A:507:HIS:HB3	2.56	0.41
1:B:91:ASP:OD1	1:D:461:GLU:OE1	2.39	0.41
1:B:148:THR:HB	1:B:279:LEU:HB3	2.03	0.41
1:D:278:ARG:HH12	1:D:487:GLU:CD	2.25	0.41
1:B:682:ASN:HB3	1:B:707:THR:HG21	2.02	0.41
1:B:669:CYS:SG	1:B:670:GLY:N	2.94	0.40
1:A:144:LEU:HD11	1:A:370:VAL:HG13	2.03	0.40
1:A:461:GLU:OE1	1:C:91:ASP:OD1	2.38	0.40
1:C:512:TRP:CH2	1:C:520:GLN:HB3	2.56	0.40
1:B:184:GLY:HA3	2:B:760:HDD:HMA2	2.04	0.40
2:A:760:HDD:HMC1	2:A:760:HDD:HBC1	2.03	0.40
1:D:511:PHE:O	1:D:515:GLN:HG2	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	725/753 (96%)	702 (97%)	19 (3%)	4 (1%)	30 17

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	725/753 (96%)	704 (97%)	17 (2%)	4 (1%)	30	17
1	C	725/753 (96%)	708 (98%)	16 (2%)	1 (0%)	56	46
1	D	725/753 (96%)	708 (98%)	13 (2%)	4 (1%)	30	17
All	All	2900/3012 (96%)	2822 (97%)	65 (2%)	13 (0%)	39	27

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	SER
1	B	725	ASP
1	D	28	SER
1	D	750	LYS
1	D	751	ILE
1	A	75	SER
1	A	709	LYS
1	A	711	ALA
1	C	75	SER
1	B	28	SER
1	B	75	SER
1	B	584	LYS
1	D	75	SER

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	612/636 (96%)	598 (98%)	14 (2%)	58	51
1	B	612/636 (96%)	586 (96%)	26 (4%)	36	24
1	C	612/636 (96%)	581 (95%)	31 (5%)	29	17
1	D	612/636 (96%)	583 (95%)	29 (5%)	32	20
All	All	2448/2544 (96%)	2348 (96%)	100 (4%)	37	25

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	LEU
1	A	73	LYS
1	A	185	PHE
1	A	205	ILE
1	A	227	TRP
1	A	237	ASP
1	A	252	ASN
1	A	375	LEU
1	A	415	TYR
1	A	440	TYR
1	A	459	ASN
1	A	552	LEU
1	A	709	LYS
1	A	712	ASP
1	B	27	ASP
1	B	37	ARG
1	B	185	PHE
1	B	198	LEU
1	B	205	ILE
1	B	227	TRP
1	B	252	ASN
1	B	265	SER
1	B	283	GLU
1	B	375	LEU
1	B	377	ARG
1	B	440	TYR
1	B	459	ASN
1	B	552	LEU
1	B	562	LEU
1	B	565	GLU
1	B	566	LEU
1	B	568	ASP
1	B	571	LEU
1	B	583	LYS
1	B	595	ASP
1	B	610	GLU
1	B	612	ARG
1	B	633	LEU
1	B	703	LYS
1	B	751	ILE
1	C	37	ARG
1	C	61	ARG
1	C	159	ILE

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Mol	Chain	Res	Type
1	C	185	PHE
1	C	205	ILE
1	C	227	TRP
1	C	237	ASP
1	C	252	ASN
1	C	375	LEU
1	C	377	ARG
1	C	440	TYR
1	C	459	ASN
1	C	478	LYS
1	C	488	ARG
1	C	521	ARG
1	C	531	LEU
1	C	552	LEU
1	C	568	ASP
1	C	571	LEU
1	C	584	LYS
1	C	606	LEU
1	C	612	ARG
1	C	616	LEU
1	C	633	LEU
1	C	648	LEU
1	C	660	LEU
1	C	685	LEU
1	C	709	LYS
1	C	725	ASP
1	C	733	LEU
1	C	750	LYS
1	D	28	SER
1	D	41	GLU
1	D	48	GLN
1	D	73	LYS
1	D	185	PHE
1	D	205	ILE
1	D	227	TRP
1	D	237	ASP
1	D	252	ASN
1	D	265	SER
1	D	340	LEU
1	D	344	GLU
1	D	375	LEU
1	D	377	ARG

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Mol	Chain	Res	Type
1	D	440	TYR
1	D	459	ASN
1	D	490	GLU
1	D	552	LEU
1	D	554	LEU
1	D	582	LEU
1	D	598	VAL
1	D	616	LEU
1	D	633	LEU
1	D	648	LEU
1	D	713	GLN
1	D	747	LYS
1	D	749	ASP
1	D	750	LYS
1	D	751	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	252	ASN
1	A	459	ASN
1	A	515	GLN
1	B	252	ASN
1	B	459	ASN
1	B	507	HIS
1	B	629	HIS
1	C	252	ASN
1	C	459	ASN
1	C	507	HIS
1	C	556	GLN
1	C	572	ASN
1	C	629	HIS
1	C	671	ASN
1	D	48	GLN
1	D	252	ASN
1	D	449	HIS
1	D	459	ASN
1	D	507	HIS
1	D	556	GLN
1	D	629	HIS
1	D	671	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	HDD	A	760	1,3	30,52,52	3.73	13 (43%)	20,89,89	3.23	10 (50%)
2	HDD	B	760	1,3	30,52,52	3.37	13 (43%)	20,89,89	3.81	13 (65%)
2	HDD	C	760	1,3	30,52,52	3.05	11 (36%)	20,89,89	3.86	13 (65%)
2	HDD	D	760	1	30,52,52	3.30	13 (43%)	20,89,89	3.68	14 (70%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HDD	A	760	1,3	-	0/3/89/89	0/1/9/9
2	HDD	B	760	1,3	-	0/3/89/89	0/1/9/9
2	HDD	C	760	1,3	-	0/3/89/89	0/1/9/9
2	HDD	D	760	1	-	0/3/89/89	0/1/9/9

All (50) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	760	HDD	C2D-C3D	-10.86	1.27	1.54
2	C	760	HDD	C2D-C3D	-10.62	1.28	1.54
2	A	760	HDD	C2D-C3D	-10.58	1.28	1.54
2	D	760	HDD	C2D-C3D	-10.38	1.28	1.54
2	B	760	HDD	C3C-C2C	-4.52	1.34	1.40
2	C	760	HDD	C3B-C2B	-4.20	1.34	1.40
2	A	760	HDD	C3B-C2B	-4.11	1.34	1.40
2	D	760	HDD	C3B-C2B	-4.00	1.35	1.40
2	C	760	HDD	C3C-C2C	-3.74	1.35	1.40
2	A	760	HDD	C3C-C2C	-3.71	1.35	1.40
2	D	760	HDD	O1D-CGD	-3.51	1.29	1.35
2	D	760	HDD	C3C-C2C	-3.51	1.35	1.40
2	B	760	HDD	C3B-C2B	-3.50	1.35	1.40
2	A	760	HDD	O1D-CGD	-3.31	1.30	1.35
2	C	760	HDD	O1D-CGD	-3.21	1.30	1.35
2	B	760	HDD	O1D-CGD	-3.19	1.30	1.35
2	B	760	HDD	C4D-ND	-2.64	1.34	1.38
2	C	760	HDD	C4D-ND	-2.45	1.34	1.38
2	D	760	HDD	C4D-ND	-2.06	1.35	1.38
2	A	760	HDD	C4B-NB	2.02	1.39	1.36
2	D	760	HDD	CMC-C2C	2.10	1.56	1.51
2	C	760	HDD	CMC-C2C	2.18	1.56	1.51
2	D	760	HDD	C3C-CAC	2.21	1.52	1.47
2	B	760	HDD	CMC-C2C	2.23	1.56	1.51
2	A	760	HDD	CMC-C2C	2.62	1.57	1.51
2	C	760	HDD	C3C-CAC	2.75	1.53	1.47
2	B	760	HDD	C3C-CAC	2.76	1.53	1.47
2	A	760	HDD	C3B-CAB	2.84	1.53	1.47
2	A	760	HDD	CMD-C2D	2.84	1.57	1.53
2	A	760	HDD	C3C-CAC	2.92	1.54	1.47
2	B	760	HDD	C3B-CAB	3.12	1.54	1.47
2	D	760	HDD	C3B-CAB	3.40	1.54	1.47
2	C	760	HDD	C3B-CAB	3.50	1.55	1.47
2	B	760	HDD	FE-ND	3.78	2.10	1.95
2	A	760	HDD	FE-ND	3.87	2.11	1.95
2	C	760	HDD	FE-ND	4.13	2.12	1.95
2	B	760	HDD	CAD-C3D	4.31	1.61	1.53
2	B	760	HDD	OND-C2D	4.46	1.51	1.42
2	A	760	HDD	OND-C2D	4.46	1.51	1.42
2	D	760	HDD	CAD-C3D	4.49	1.61	1.53
2	D	760	HDD	FE-ND	4.71	2.14	1.95
2	C	760	HDD	OND-C2D	4.81	1.52	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	760	HDD	OND-C2D	4.87	1.52	1.42
2	A	760	HDD	O1D-C3D	5.14	1.55	1.46
2	D	760	HDD	O1D-C3D	5.54	1.56	1.46
2	C	760	HDD	O1D-C3D	5.73	1.56	1.46
2	D	760	HDD	CMD-C2D	5.78	1.61	1.53
2	B	760	HDD	O1D-C3D	6.12	1.57	1.46
2	B	760	HDD	CMD-C2D	6.15	1.62	1.53
2	A	760	HDD	CAD-C3D	12.03	1.76	1.53

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	760	HDD	OND-C2D-CMD	-10.11	91.16	109.41
2	D	760	HDD	OND-C2D-CMD	-8.13	94.74	109.41
2	B	760	HDD	OND-C2D-CMD	-7.26	96.29	109.41
2	B	760	HDD	CAA-C2A-C1A	-6.74	119.69	127.01
2	D	760	HDD	CAA-C2A-C1A	-6.23	120.25	127.01
2	B	760	HDD	CAA-CBA-CGA	-6.17	101.44	112.75
2	C	760	HDD	C3B-C4B-NB	-5.78	100.03	110.94
2	B	760	HDD	C3B-C4B-NB	-5.72	100.14	110.94
2	A	760	HDD	C3B-C4B-NB	-5.65	100.28	110.94
2	D	760	HDD	C3B-C4B-NB	-5.36	100.82	110.94
2	D	760	HDD	CAA-CBA-CGA	-5.28	103.07	112.75
2	A	760	HDD	C3C-CAC-CBC	-5.25	115.57	126.32
2	A	760	HDD	O1D-CGD-CBD	-5.13	104.19	110.20
2	B	760	HDD	C2D-C1D-CHD	-5.09	115.56	123.48
2	C	760	HDD	O1D-CGD-CBD	-4.90	104.46	110.20
2	A	760	HDD	OND-C2D-CMD	-4.87	100.61	109.41
2	C	760	HDD	C3C-CAC-CBC	-4.74	116.61	126.32
2	A	760	HDD	CAA-CBA-CGA	-4.61	104.29	112.75
2	D	760	HDD	C3C-C4C-NC	-4.49	103.41	109.21
2	A	760	HDD	C3C-C4C-NC	-4.41	103.51	109.21
2	A	760	HDD	CAA-C2A-C1A	-4.34	122.30	127.01
2	C	760	HDD	CAA-CBA-CGA	-4.26	104.93	112.75
2	C	760	HDD	C3C-C4C-NC	-4.18	103.81	109.21
2	C	760	HDD	C2D-C1D-CHD	-4.10	117.09	123.48
2	B	760	HDD	C3C-CAC-CBC	-3.98	118.17	126.32
2	B	760	HDD	C3C-C4C-NC	-3.72	104.40	109.21
2	D	760	HDD	C3C-CAC-CBC	-3.63	118.90	126.32
2	B	760	HDD	O1D-CGD-CBD	-3.55	106.05	110.20
2	D	760	HDD	C2D-C1D-CHD	-3.48	118.06	123.48
2	C	760	HDD	C3B-CAB-CBB	-3.35	119.46	126.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	760	HDD	C2D-C1D-CHD	-3.05	118.72	123.48
2	B	760	HDD	C3B-CAB-CBB	-3.05	120.08	126.32
2	D	760	HDD	C3B-CAB-CBB	-2.97	120.24	126.32
2	D	760	HDD	O1D-CGD-CBD	-2.88	106.83	110.20
2	C	760	HDD	CAA-C2A-C1A	-2.84	123.92	127.01
2	A	760	HDD	CMA-C3A-C4A	-2.26	124.62	128.36
2	B	760	HDD	CMC-C2C-C1C	-2.26	124.62	128.36
2	C	760	HDD	CAA-C2A-C3A	-2.11	122.99	129.00
2	C	760	HDD	O1D-CGD-O2D	2.01	122.71	120.80
2	B	760	HDD	O1D-CGD-O2D	2.30	122.99	120.80
2	B	760	HDD	CMA-C3A-C2A	2.38	130.21	125.24
2	D	760	HDD	CBA-CAA-C2A	2.39	116.81	112.53
2	D	760	HDD	CMA-C3A-C2A	2.42	130.30	125.24
2	C	760	HDD	CMA-C3A-C2A	2.46	130.38	125.24
2	D	760	HDD	CMB-C2B-C3B	2.49	129.96	125.09
2	A	760	HDD	CMA-C3A-C2A	2.86	131.22	125.24
2	D	760	HDD	O1D-CGD-O2D	2.91	123.57	120.80
2	D	760	HDD	C4D-ND-C1D	4.00	109.94	107.36
2	B	760	HDD	C4D-ND-C1D	4.19	110.07	107.36
2	C	760	HDD	C4D-ND-C1D	4.39	110.19	107.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	760	HDD	6	0
2	B	760	HDD	3	0
2	C	760	HDD	2	0
2	D	760	HDD	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	727/753 (96%)	-0.45	12 (1%) 73 76	3, 8, 24, 47	1 (0%)
1	B	727/753 (96%)	-0.38	17 (2%) 64 67	4, 9, 27, 45	1 (0%)
1	C	727/753 (96%)	-0.39	17 (2%) 64 67	3, 9, 26, 44	1 (0%)
1	D	727/753 (96%)	-0.46	9 (1%) 81 83	3, 9, 25, 45	1 (0%)
All	All	2908/3012 (96%)	-0.42	55 (1%) 70 73	3, 9, 26, 47	4 (0%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	27	ASP	9.4
1	A	711	ALA	8.9
1	B	27	ASP	6.3
1	B	726	GLY	6.3
1	C	28	SER	6.2
1	A	28	SER	6.1
1	A	710	ILE	5.8
1	C	27	ASP	5.5
1	D	27	ASP	5.4
1	A	712	ASP	5.1
1	A	713	GLN	5.0
1	A	32	GLU	4.4
1	D	749	ASP	4.1
1	D	32	GLU	3.7
1	B	750	LYS	3.7
1	B	32	GLU	3.6
1	D	28	SER	3.2
1	C	594	PRO	3.2
1	B	713	GLN	3.1
1	C	749	ASP	3.1
1	B	612	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	751	ILE	3.0
1	C	726	GLY	3.0
1	B	28	SER	3.0
1	D	750	LYS	2.7
1	A	596	GLY	2.7
1	C	703	LYS	2.7
1	A	749	ASP	2.6
1	C	32	GLU	2.6
1	C	725	ASP	2.6
1	C	751	ILE	2.6
1	A	750	LYS	2.5
1	B	749	ASP	2.5
1	B	677	ASP	2.5
1	A	709	LYS	2.5
1	D	583	LYS	2.4
1	C	750	LYS	2.4
1	A	617	LEU	2.4
1	C	617	LEU	2.3
1	B	712	ASP	2.2
1	C	645	GLY	2.2
1	B	568	ASP	2.2
1	B	583	LYS	2.2
1	C	572	ASN	2.2
1	C	595	ASP	2.2
1	B	572	ASN	2.2
1	D	33	ASP	2.2
1	B	725	ASP	2.2
1	C	568	ASP	2.1
1	B	673	ALA	2.1
1	B	37	ARG	2.1
1	D	612	ARG	2.1
1	B	594	PRO	2.0
1	C	646	THR	2.0
1	C	675	ILE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	HDD	B	760	44/44	0.98	0.07	0.10	4,6,9,11	0
2	HDD	D	760	44/44	0.98	0.07	-0.00	3,5,9,10	0
2	HDD	A	760	44/44	0.98	0.07	0.00	3,5,9,10	0
2	HDD	C	760	44/44	0.98	0.07	-0.16	4,7,11,13	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.